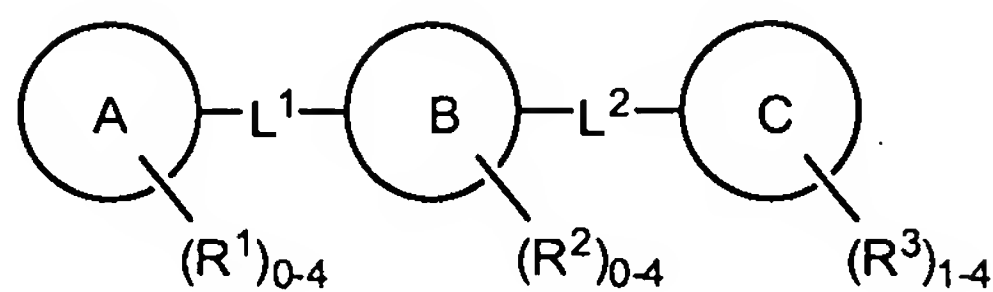


The Listing of the Claims:

1. (currently amended) A compound for modulating c-Kit activity according to Formula I,



I

or a pharmaceutically acceptable salt, thereof, wherein,

ring A is:


wherein each Y is independently either =C(H)- or =N-; and Z is selected from -O-, -S-, and -N(R<sup>7</sup>)-;

each R<sup>1</sup> is independently selected from -H, halogen, trihalomethyl, -CN, -NO<sub>2</sub>, -OR<sup>4</sup>, -S(O)<sub>0-2</sub>R<sup>4</sup>, -SO<sub>2</sub>N(R<sup>4</sup>)R<sup>4</sup>, -C(=O)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)R<sup>4</sup>, -N(R<sup>4</sup>)SO<sub>2</sub>R<sup>4</sup>, -N(R<sup>4</sup>)C(O)R<sup>4</sup>, -NCO<sub>2</sub>R<sup>4</sup>, -C(=O)R<sup>4</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

two adjacent of R<sup>1</sup>, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R<sup>10</sup>;

L<sup>1</sup> is a single bond;

ring B is phenyl ~~a five to ten membered aryl or a five to ten membered heterocyclyl~~;

each R<sup>2</sup> is independently selected from -H, halogen, trihalomethyl, -CN, -NO<sub>2</sub>, -OR<sup>4</sup>, -N(R<sup>4</sup>)R<sup>4</sup>, -S(O)<sub>0-2</sub>R<sup>4</sup>, -SO<sub>2</sub>N(R<sup>4</sup>)R<sup>4</sup>, -C(=O)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)R<sup>4</sup>, -N(R<sup>4</sup>)SO<sub>2</sub>R<sup>4</sup>, -N(R<sup>4</sup>)C(O)R<sup>4</sup>, -NCO<sub>2</sub>R<sup>4</sup>, -C(=O)R<sup>4</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

two adjacent of R<sup>2</sup>, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R<sup>15</sup>;

L<sup>2</sup> is selected from -N(H)N(H)C(=O)N(H)-, -CH<sub>2</sub>N(H)C(=O)N(H)-, -CH<sub>2</sub>OC(=O)N(H)-, and -XCH<sub>2</sub>C(=O)N(H)-; wherein X is selected from -O-, -S(O)<sub>0-2</sub>-, and -N(R<sup>7</sup>)-; and any C-H of L<sup>2</sup> is optionally C-R<sup>20</sup>;

ring C is phenyl ~~or pyridyl~~;

each R<sup>3</sup> is independently selected from halogen, trihalomethyl, -CN,

-NO<sub>2</sub>, -OR<sup>4</sup>, -N(R<sup>4</sup>)R<sup>4</sup>, -S(O)<sub>0-2</sub>R<sup>4</sup>, -SO<sub>2</sub>N(R<sup>4</sup>)R<sup>4</sup>, -C(=O)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)R<sup>4</sup>, -N(R<sup>4</sup>)SO<sub>2</sub>R<sup>4</sup>, -N(R<sup>4</sup>)C(O)R<sup>4</sup>, -NCO<sub>2</sub>R<sup>4</sup>, -C(=O)R<sup>4</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl; provided R<sup>3</sup> is not a cyclic sulfonamide attached to ring C via the nitrogen of said cyclic sulfonamide,

wherein there exists at least one of R<sup>3</sup> that is halogen or trihalomethyl;

~~two adjacent of R<sup>3</sup>, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R<sup>25</sup>;~~

R<sup>4</sup> is selected from -H, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

two of R<sup>4</sup>, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R<sup>5</sup> is selected from -H, -CN, -NO<sub>2</sub>, -OR<sup>4</sup>, -S(O)<sub>0-2</sub>R<sup>4</sup>, -CO<sub>2</sub>R<sup>4</sup>, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted C<sub>1-6</sub>alkenyl, and optionally substituted C<sub>1-6</sub>alkynyl;

R<sup>7</sup> is selected from -H, optionally substituted C<sub>1-6</sub>alkyl, -SO<sub>2</sub>N(R<sup>4</sup>)R<sup>4</sup>, -CO<sub>2</sub>R<sup>4</sup>, -C(=O)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)R<sup>4</sup>, -C(=O)R<sup>4</sup>, optionally substituted alkoxy, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl; and

each of R<sup>10</sup>, each of R<sup>15</sup>, each of R<sup>20</sup>, and each of R<sup>25</sup> is independently selected from -H, halogen, trihalomethyl, -CN, -NO<sub>2</sub>, -OR<sup>4</sup>, -N(R<sup>4</sup>)R<sup>4</sup>, -S(O)<sub>0-2</sub>R<sup>4</sup>, -SO<sub>2</sub>N(R<sup>4</sup>)R<sup>4</sup>, -CO<sub>2</sub>R<sup>4</sup>, -C(=O)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)R<sup>4</sup>, -N(R<sup>4</sup>)SO<sub>2</sub>R<sup>4</sup>, -N(R<sup>4</sup>)C(O)R<sup>4</sup>, -NCO<sub>2</sub>R<sup>4</sup>, -C(=O)R<sup>4</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl,

optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

provided:

1) ~~when both ring B and ring C are phenyl:~~

a) ~~and the compound comprises ring B-CH<sub>2</sub>N(H)C(=O)N(H)-ring C, then L<sup>1</sup> must be a single bond; R<sup>3</sup> can not comprise a group of the formula -O(CH<sub>2</sub>)<sub>2-4</sub>-N-piperazine that is *ortho* to L<sup>2</sup>; and ring A cannot be a 5-methyl-[1,2,4]-oxadiazol-3-yl radical, a 4H-[1,2,4]-oxadiazol-5-one-3-yl radical, nor a 4'-[2,2';6',2'']terpyridinyl radical;~~

b) ~~and L<sup>1</sup> is single bond, then L<sup>2</sup> cannot comprise N(H)C(=O)C(=O)N(H) nor N(H)C(=Q)C(H)CNC(=O) (where Q is S or O);~~

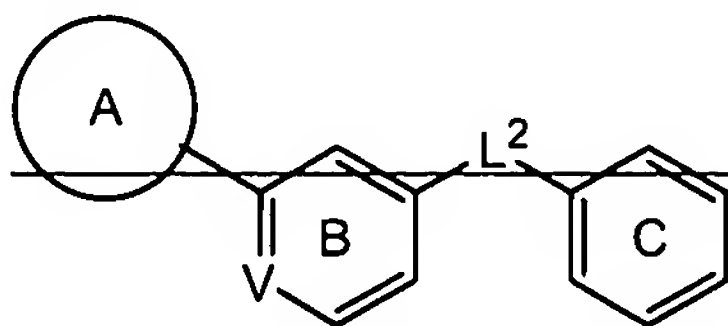
c) ~~and L<sup>1</sup> is other than single bond, then A cannot be quinolin-2-yl L<sup>1</sup>, quinolin-3-yl L<sup>1</sup>, or quinolin-4-yl L<sup>1</sup>;~~

2) ~~when ring A is a fused aryl system, then L<sup>1</sup> must be a single bond;~~

3) ~~when ring B is phenyl, ring C is a C<sub>6-16</sub>carbocyclic, L<sup>1</sup> is a single bond, and the compound comprises ring B-OCH<sub>2</sub>C(=O)N(H) then ring A cannot be a 2,5-dimethyl-1H-pyrrole-1-yl radical;~~

4) ~~ring A cannot be a pyrimidin-2-yl radical when L<sup>1</sup> is N(H) and ring B is phenyl;~~

5) ~~when the compound comprises the formula,~~



~~where V is -C(H)- or -N-, and there is a nitrogen of L<sup>2</sup> bound directly to ring B, then A can not comprise a [1,2,4]-oxadiazol-3-yl radical; and~~

6) ~~the compound is not one of: N-naphthalen-1-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-[4-(phenyloxy)phenyl]-2-[[3-(1H-tetrazol-1-~~

~~yl)phenyl]oxy}acetamide, N (3,4 dimethylphenyl) 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N (2,3 dimethylphenyl) 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N (2,4 dimethylphenyl) 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N (2,5 dimethylphenyl) 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N (3,5 dimethylphenyl) 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N (2,6 dimethylphenyl) 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy} N (2,4,6-trimethylphenyl)acetamide, N (2-ethylphenyl) 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N (4-ethylphenyl) 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N (2,6-diethylphenyl) 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N [2-(methyloxy)phenyl] 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N [2-(ethyloxy)phenyl] 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N [3-(ethyloxy)phenyl] 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N [2,4-bis(methyloxy)phenyl] 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N [4-(dimethylamino)phenyl] 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-chloro-3-methylphenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-fluorophenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2- {[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl]acetamide, 2- {[3-(1H-tetrazol-1-yl)phenyl] oxy}-N-[3-(trifluoromethyl)phenyl]acetamide, methyl 4 [({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoate, ethyl 4 [({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoate, 3 [({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoic acid, N [3-(methyloxy)phenyl] 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N [4-(methyloxy)phenyl] 2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-chloro-5-(trifluoromethyl) phenyl]-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-~~

{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide or ~~acetamide~~, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, ~~N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ and ~~N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide.~~

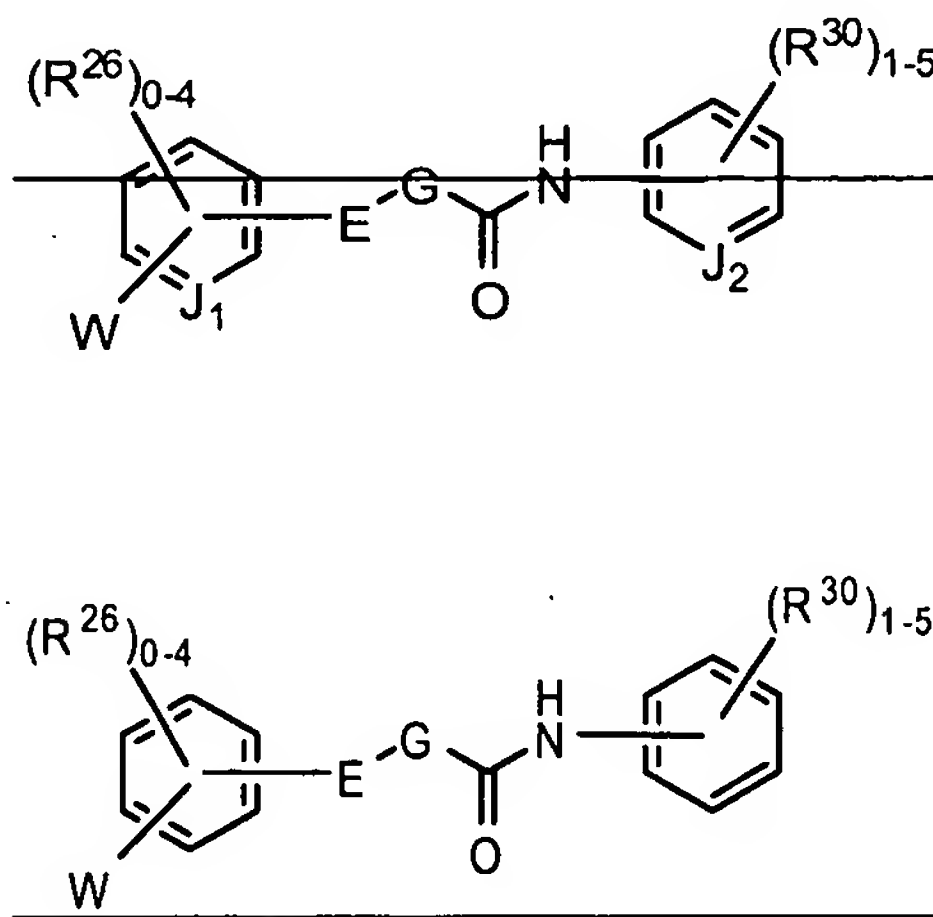
2-12. (cancelled)

13. (Currently Amended) The compound according to claim ~~10~~, 1, wherein there exists at least one of R<sup>3</sup> that is trifluoromethyl.

14. (original) The compound according to claim 13, wherein ring C is a phenyl comprising a trifluoromethyl radical *meta*- to L<sup>2</sup>.

15. (currently amended) The compound according to claim 1, ~~claim 10~~, wherein each of R<sup>3</sup> is independently selected from halogen, trihalomethyl, -OR<sup>4</sup>, -C(=O)R<sup>4</sup>, and optionally substituted C<sub>1-6</sub>alkyl.

16. (currently amended) A compound for modulating c-Kit activity according to the following Formula: H,



or a pharmaceutically acceptable salt, thereof, wherein,

W is selected from the following:


each of  $R^{27}$  independently selected from halogen, trihalomethyl,  $-\text{CN}$ ,  $-\text{NO}_2$ ,  $-\text{OR}^{55}$ ,  $-\text{S}(\text{O})_{0-2}\text{R}^{55}$ ,  $-\text{SO}_2\text{N}(\text{R}^{55})\text{R}^{55}$ ,  $-\text{C}(=\text{O})\text{N}(\text{R}^{55})\text{R}^{55}$ ,  $-\text{C}(=\text{NR}^{50})\text{N}(\text{R}^{55})\text{R}^{55}$ ,  $-\text{C}(=\text{NR}^{50})\text{R}^{55}$ ,  $-\text{N}(\text{R}^{55})\text{SO}_2\text{R}^{55}$ ,  $-\text{N}(\text{R}^{55})\text{C}(\text{O})\text{R}^{55}$ ,  $-\text{NCO}_2\text{R}^{55}$ ,  $-\text{C}(=\text{O})\text{R}^{55}$ , optionally substituted alkoxy, optionally substituted  $\text{C}_{1-6}$ alkyl, optionally substituted aryl  $\text{C}_{1-6}$ alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl  $\text{C}_{1-6}$ alkyl;

each Y is independently either  $=\text{C}(\text{H})-$  or  $=\text{N}-$ ;

Z is selected from  $-\text{O}-$ ,  $-\text{S}(\text{O})_{0-2}-$ , and  $-\text{N}(\text{R}^7)-$

E and G are each independently selected from  $-\text{O}-$ ,  $-\text{S}(\text{O})_{0-2}-$ ,  $-\text{C}(\text{R}^{31})\text{R}^{32}-$ , and  $-\text{N}(\text{R}^{33})-$ ;

$\text{J}_1$  and  $\text{J}_2$  are each independently  $=\text{C}(\text{H})-$  or  $=\text{N}-$ ;

$R^{26}$  is independently selected from -H, halogen, trihalomethyl, -CN, -NO<sub>2</sub>, -OR<sup>40</sup>, -N(R<sup>40</sup>)R<sup>40</sup>, -S(O)<sub>0-2</sub>R<sup>40</sup>, -SO<sub>2</sub>N(R<sup>40</sup>)R<sup>40</sup>, -CO<sub>2</sub>R<sup>40</sup>, -C(=O)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)R<sup>40</sup>, -N(R<sup>40</sup>)SO<sub>2</sub>R<sup>40</sup>, -N(R<sup>40</sup>)C(O)R<sup>40</sup>, -NCO<sub>2</sub>R<sup>40</sup>, -C(=O)R<sup>40</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

$R^{30}$  is independently selected from halogen, trihalomethyl, -CN, -NO<sub>2</sub>, -OR<sup>40</sup>, -N(R<sup>40</sup>)R<sup>40</sup>, -S(O)<sub>0-2</sub>R<sup>40</sup>, -SO<sub>2</sub>N(R<sup>40</sup>)R<sup>40</sup>, -C(=O)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)R<sup>40</sup>, -N(R<sup>40</sup>)SO<sub>2</sub>R<sup>40</sup>, -N(R<sup>40</sup>)C(O)R<sup>40</sup>, -NCO<sub>2</sub>R<sup>40</sup>, -C(=O)R<sup>40</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl, wherein there exists at least one of  $R^{30}$  that is trihalomethyl; or

~~two adjacent of  $R^{26}$  or two adjacent of  $R^{30}$ , together with the annular atoms to which they are attached, can form a five to six membered ring containing up to two heteroatoms and optionally substituted with up to three of  $R^{35}$ ;~~

$R^{31}$  and  $R^{32}$  are each independently selected from -H, halogen, trihalomethyl, -CN, -NO<sub>2</sub>, -OR<sup>40</sup>, -N(R<sup>40</sup>)R<sup>40</sup>, -S(O)<sub>0-2</sub>R<sup>40</sup>, -SO<sub>2</sub>N(R<sup>40</sup>)R<sup>40</sup>, -CO<sub>2</sub>R<sup>40</sup>, -C(=O)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)R<sup>40</sup>, -N(R<sup>40</sup>)SO<sub>2</sub>R<sup>40</sup>, -N(R<sup>40</sup>)C(O)R<sup>40</sup>, -NCO<sub>2</sub>R<sup>40</sup>, -C(=O)R<sup>40</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

$R^{33}$  is selected from -H, optionally substituted lower alkyl, -SO<sub>2</sub>N(R<sup>40</sup>)R<sup>40</sup>, -CO<sub>2</sub>R<sup>40</sup>, -C(=O)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)R<sup>40</sup>, -C(=O)R<sup>40</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

$R^{40}$  is selected from -H, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl,

and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

two of R<sup>40</sup>, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R<sup>50</sup> is selected from -H, -CN, -NO<sub>2</sub>, -OR<sup>40</sup>, -S(O)<sub>0-2</sub>R<sup>40</sup>, -CO<sub>2</sub>R<sup>40</sup>, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted C<sub>1-6</sub>alkenyl, and optionally substituted C<sub>1-6</sub>alkynyl;

R<sup>55</sup> is selected from -H, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl; and

two of R<sup>55</sup>, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P.

17. (cancelled)

18. (currently amended) The compound according to claim ~~17~~ 16, wherein R<sup>30</sup> is selected from halogen, trihalomethyl, -OR<sup>40</sup>, -N(R<sup>40</sup>)R<sup>40</sup>, -C(=O)R<sup>40</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl, wherein there exists at least one of R<sup>30</sup> that is trifluoromethyl.

19. (cancelled)

20. (cancelled)

21. (cancelled)

22. (cancelled)

23. (withdrawn from consideration, currently amended) The compound according to ~~claim 22~~, claim 16, wherein E is selected from -O-, -S(O)<sub>0-2</sub>-, and -NH-; and G is -CH<sub>2</sub>-.

24. (withdrawn from consideration, currently amended) The compound according to ~~claim 22~~, claim 16, wherein E is either -CH<sub>2</sub>- or -NH-; and G is selected from -O-, -S-, and -NH-.

25. (cancelled)

26. (cancelled)

27. (currently amended) ~~The compound according to claim 1, selected from Table 3:~~

A compound selected from the following Table:

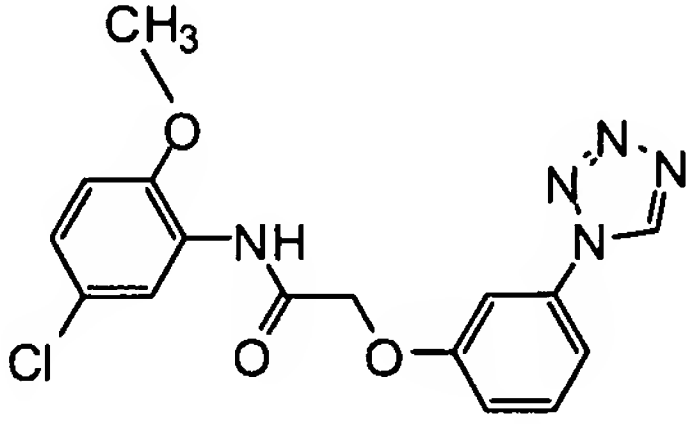
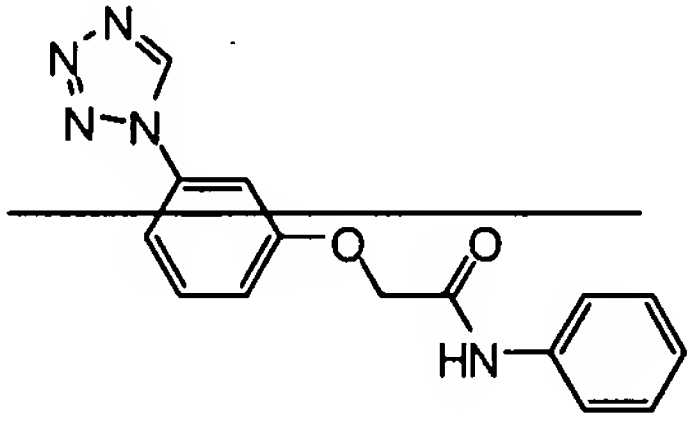
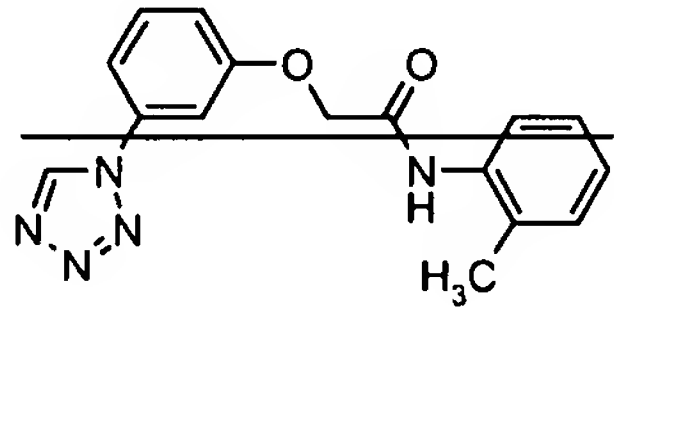
Table 3		
Entry	Name	Structure
1	N-[5-chloro-2-(methoxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
2	<del>N-phenyl 2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide</del>	
3	<del>N-(2-methylphenyl) 2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide</del>	

Table 3		
Entry	Name	Structure
4	N-(2-chlorophenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
5	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
6	<del>ethyl 2- [({[3-(1H-tetrazol-1-yl)phenyl]oxy} acetyl)amino]-4,5,6,7-tetrahydro-1-benzothiophene-3-carboxylate</del>	
7	N-(3-chloro-2-methylphenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
8	N-(3-fluorophenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	

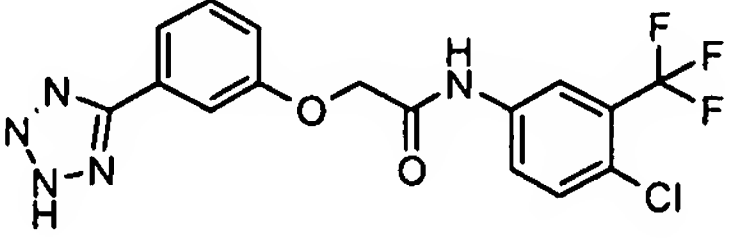
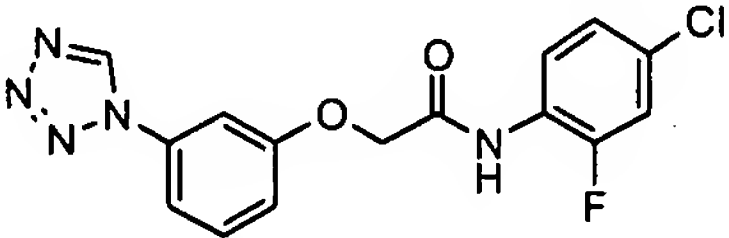
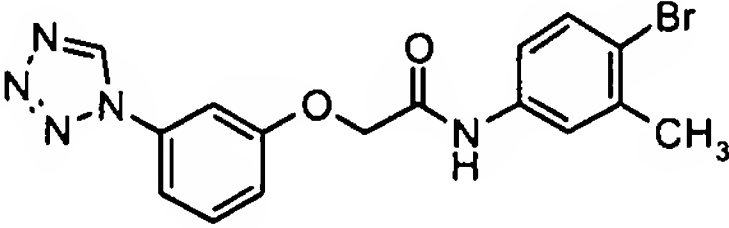
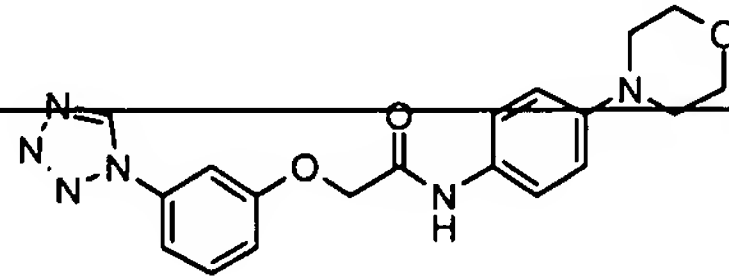
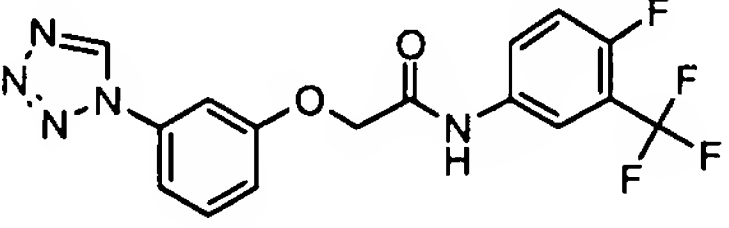
Table 3		
Entry	Name	Structure
9	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(2H-tetrazol-5- yl)phenyl]oxy}acetamide	
10	N-(4-chloro-2-fluorophenyl)-2- {[3-(1H- tetrazol-1-yl)phenyl]oxy}acetamide	
11	N-(4-bromo-3-methylphenyl)-2- {[3-(1H- tetrazol-1-yl)phenyl]oxy}acetamide	
12	<del>N-(4-morpholin-4-ylphenyl)-2- {[3-(1H- tetrazol-1-yl)phenyl]oxy}acetamide</del>	
13	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	

Table 3		
Entry	Name	Structure
14	N-[4-bromo-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	
15	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[4-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	
16	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]oxy}propanamide	
17	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(5-methyl-1H-tetrazol-1- yl)phenyl]oxy}acetamide	
18	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-methyl-5-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	

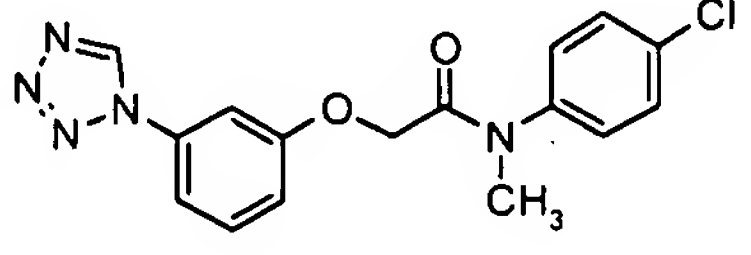
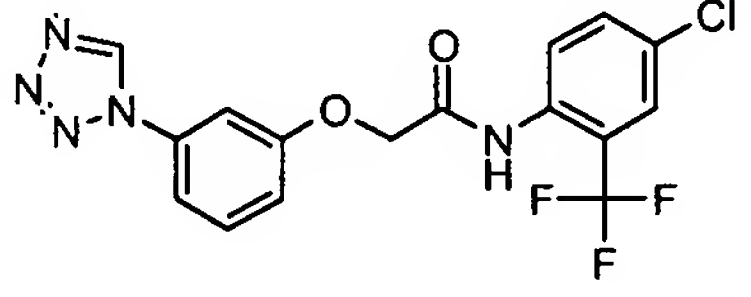
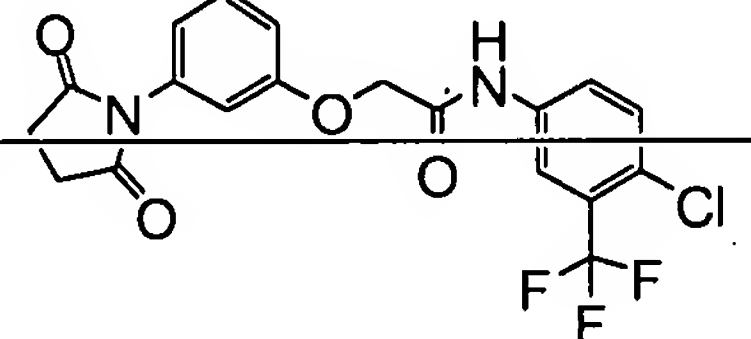
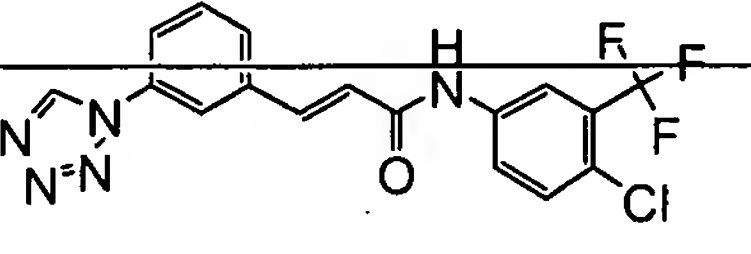
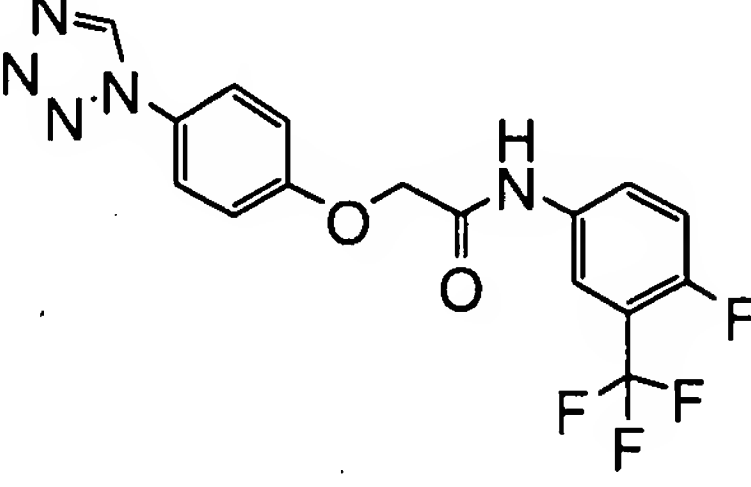
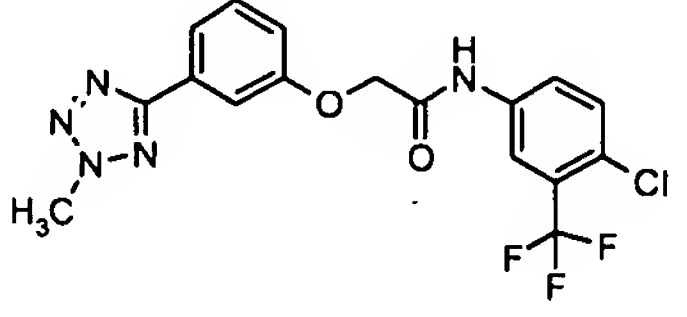
Table 3		
Entry	Name	Structure
19	N-(4-chlorophenyl)-N-methyl-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
20	N-[4-chloro-2-(trifluoromethyl)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
21	<del>N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-(2,5-dioxopyrrolidin-1-yl)phenyl]oxy}acetamide</del>	
22	<del>(2E)-N-[4-chloro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]prop-2-enamide</del>	
23	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
24	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-(2-methyl-2H-tetrazol-5-yl)phenyl]oxy}acetamide	

Table 3		
Entry	Name	Structure
25	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2,4-dichloro-5-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	
26	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]thio}acetamide	
27	N-[4-chloro-3-(trifluoromethyl)phenyl]- N~2~-[3-(1H-tetrazol-1- yl)phenyl]glycinamide	
28	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	
29	<del>methyl 1-{3-[(2-{[4-chloro-3- (trifluoromethyl)phenyl]amino}-2- oxoethyl)oxy]phenyl}-1H-1,2,3-triazole-4- carboxylate</del>	

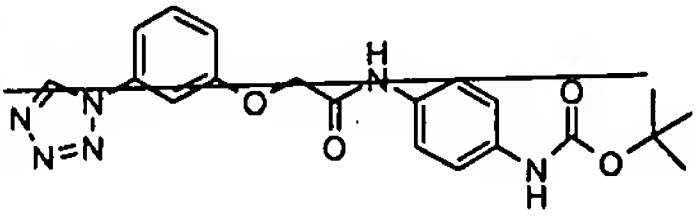
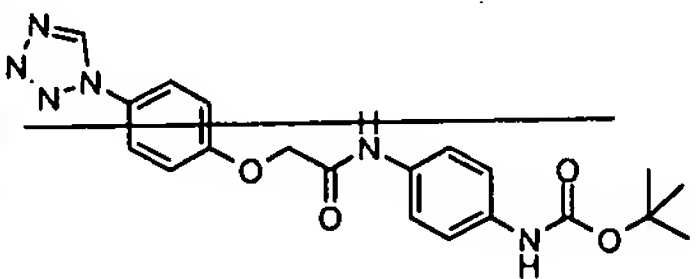
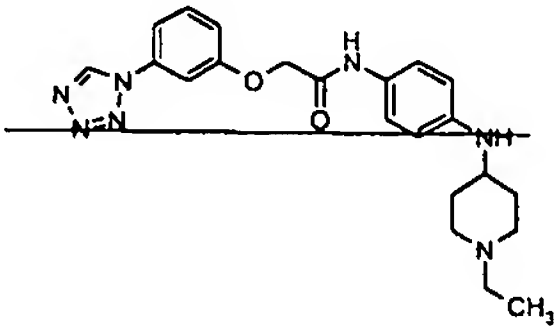
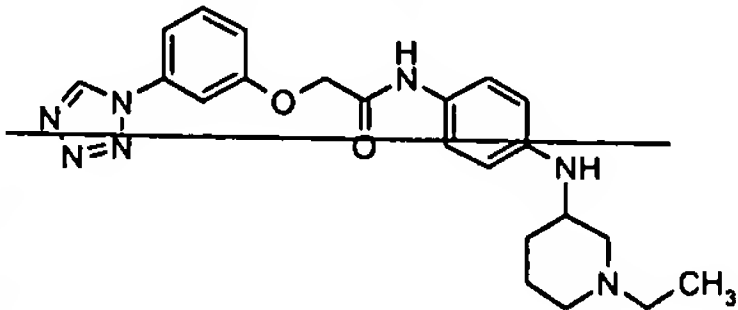
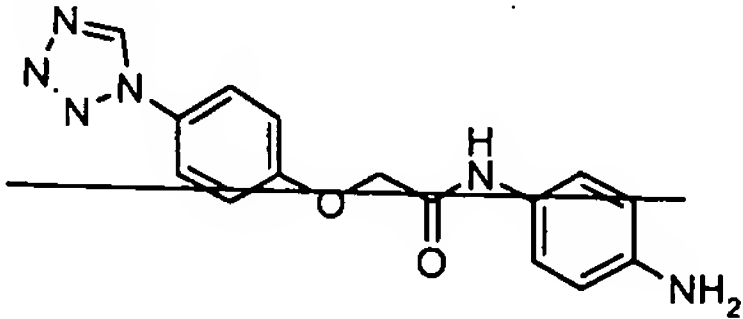
Table 3		
Entry	Name	Structure
30	1,1 dimethylethyl {4 [({[3 (1H tetrazol-1-yl)phenyl]oxy}acetyl)amino]phenyl}carbamate	
31	1,1 dimethylethyl {4 [({[4 (1H tetrazol-1-yl)phenyl]oxy}acetyl)amino]phenyl}carbamate	
32	N {4 [(1-ethylpiperidin-4-yl)amino]phenyl} 2 {[3 (1H tetrazol-1-yl)phenyl]oxy}acetamide	
33	N {4 [(1-ethylpiperidin-3-yl)amino]phenyl} 2 {[3 (1H tetrazol-1-yl)phenyl]oxy}acetamide	
34	N (4-aminophenyl) 2 {[4 (1H tetrazol-1-yl)phenyl]oxy}acetamide	

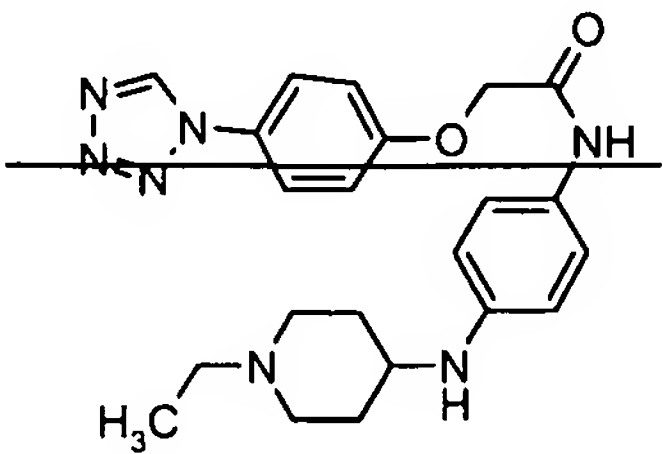
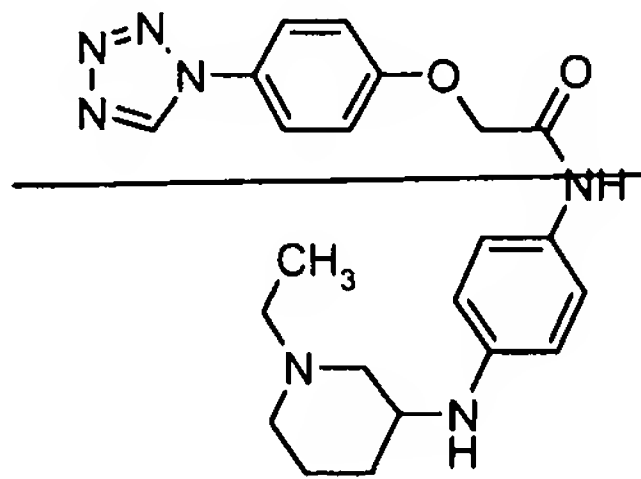
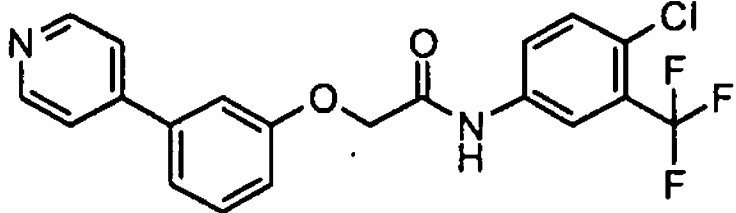
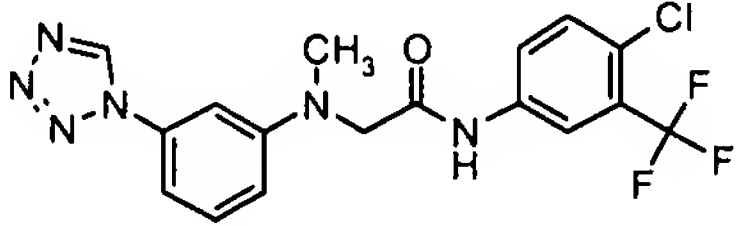
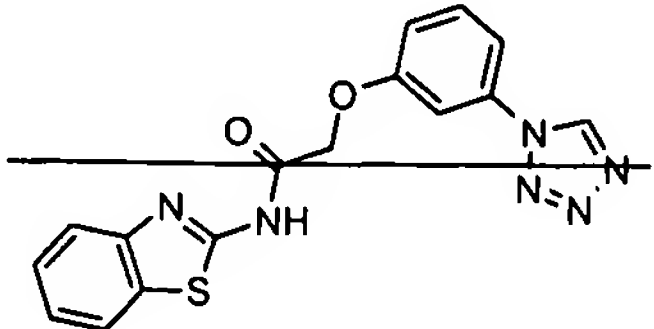
Table 3		
Entry	Name	Structure
35	N-{4-[(1-ethylpiperidin-4-yl)amino]phenyl}-2-[[4-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
36	N-{4-[(1-ethylpiperidin-3-yl)amino]phenyl}-2-[[4-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
37	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyridin-4-ylphenyl)oxy]acetamide	
38	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-methyl-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
39	N-1,3-benzothiazol-2-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	

Table 3		
Entry	Name	Structure
40	N-quinolin-8-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
41	N-(2,3-dihydro-1,4-benzodioxin-6-yl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
42	N-isoquinolin-5-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
43	N-{3-[(phenylmethyl)oxy]phenyl}-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
44	N-[5-methyl-2-(methoxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	

Table 3		
Entry	Name	Structure
45	N-[2,5-bis(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
46	N-(6-fluoro-1,3-benzothiazol-2-yl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
47	methyl 3-(((3-(1H-tetrazol-1-yl)phenyl)oxy)acetyl)amino]benzoate	
48	5-chloro-2-(((3-(1H-tetrazol-1-yl)phenyl)oxy)acetyl)amino]benzamide	
49	N-[5-chloro-2,4-bis(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	

Table 3		
Entry	Name	Structure
50	N-[2-(phenyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
51	N-[3-(aminosulfonyl)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
52	N-[2-(methyloxy)-5-(trifluoromethyl)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
53	N-(4-[[4-methylphenyl)sulfonyl]amino]phenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
54	N-(5-phenyl-1H-pyrazol-3-yl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	

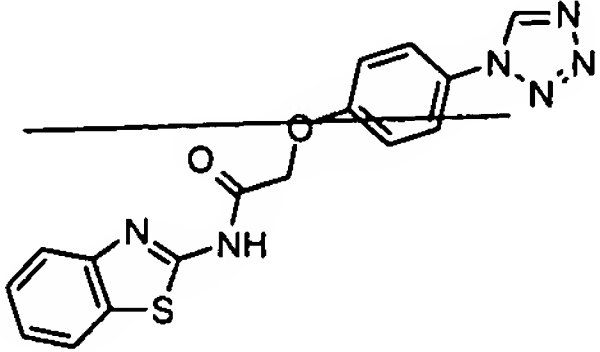
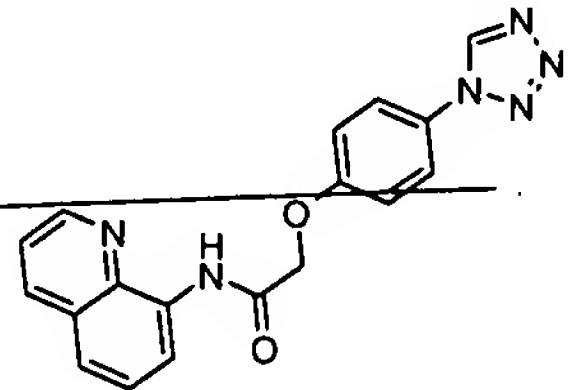
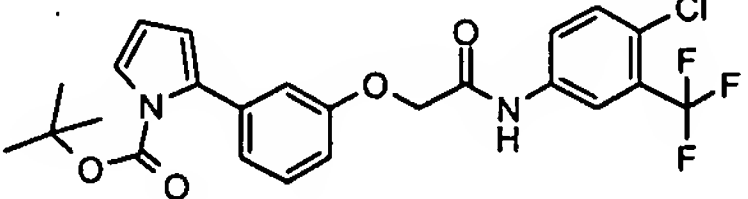
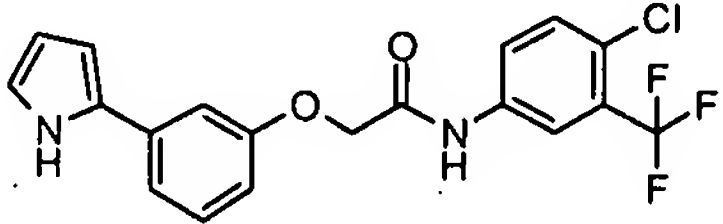
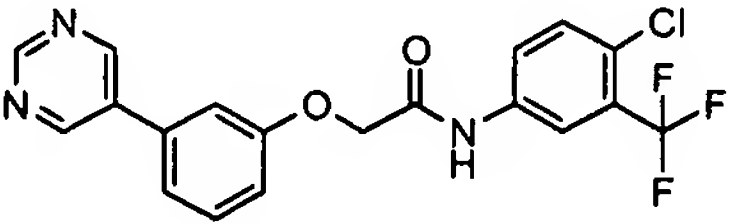
Table 3		
Entry	Name	Structure
55	<del>N-1,3-benzothiazol-2-yl-2-[[4-(1H-tetrazol-1-yl)phenyl]oxy]acetamide</del>	
56	<del>N-quinolin-8-yl-2-[[4-(1H-tetrazol-1-yl)phenyl]oxy]acetamide</del>	
57	1,1-dimethylethyl 2-{3-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-pyrrole-1-carboxylate	
58	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-pyrrol-2-yl)phenyl]oxy}acetamide	
59	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyrimidin-5-ylphenyl)oxy]acetamide	

Table 3		
Entry	Name	Structure
60	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-1,2,3-triazol-1- yl)phenyl]oxy}acetamide	
61	<del>4-chloro N (2 {[3 (1H tetrazol 1- yl)phenyl]oxy}ethyl) 3- (trifluoromethyl)aniline</del>	
62	<del>N [4 chloro 3 (trifluoromethyl)phenyl] N- (2 {[3 (1H tetrazol 1- yl)phenyl]oxy}ethyl)formamide</del>	
63	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-pyridin-3-ylphenyl)oxy]acetamide	
64	<del>N [4 chloro 3 (trifluoromethyl)phenyl] 2- [(3 furan 3 ylphenyl)oxy]acetamide</del>	
65	<del>(2E) N [4 fluoro 3- (trifluoromethyl)phenyl] 3 [3 (1H- tetrazol 1 yl)phenyl]prop 2 enamide</del>	

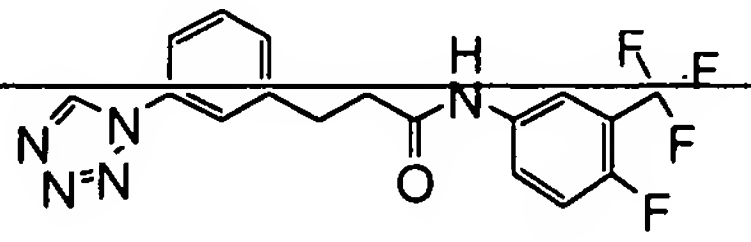
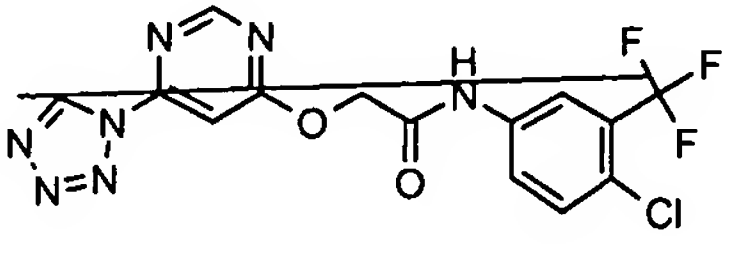
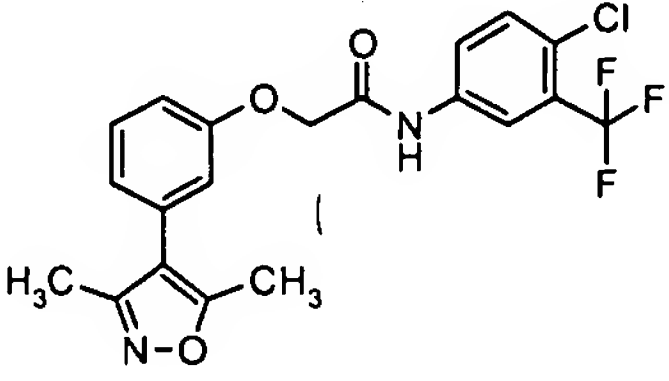
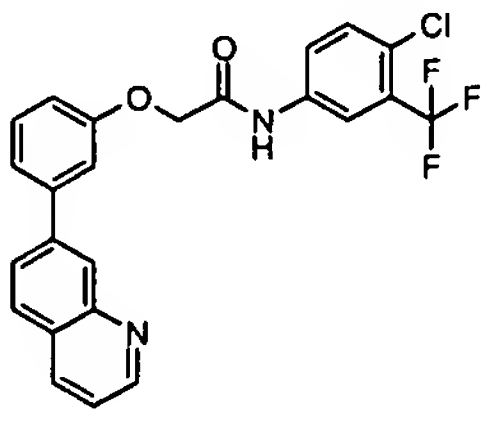
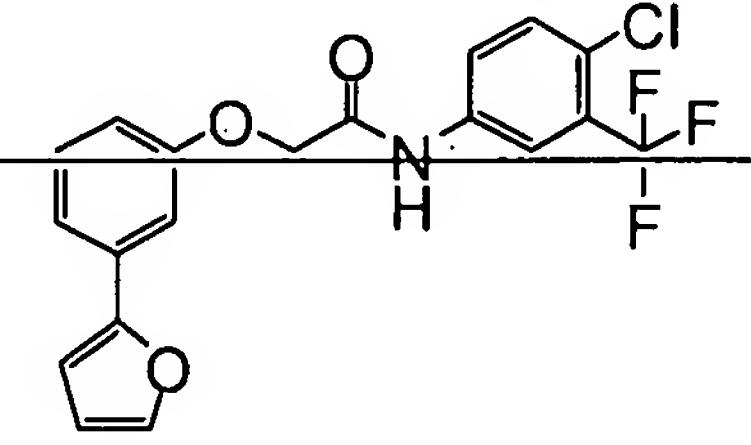
Table 3		
Entry	Name	Structure
66	<del>N-[4-fluoro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]propanamide</del>	
67	<del>N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[6-(1H-tetrazol-1-yl)pyrimidin-4-yl]oxy]acetamide</del>	
68	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[3-(3,5-dimethylisoxazol-4-yl)phenyl]oxy]acetamide	
69	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[3-quinolin-7-ylphenyl]oxy]acetamide	
70	<del>N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[3-furan-2-ylphenyl]oxy]acetamide</del>	

Table 3		
Entry	Name	Structure
71	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
72	<del>N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-dibenzo[b,d]furan-4-ylphenyl)oxy]acetamide</del>	
73	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
74	<del>N-methyl N-[4-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide</del>	
75	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[[3-(1H-tetrazol-1-yl)phenyl]methyl]urea	

Table 3		
Entry	Name	Structure
76	<del>N-[4-chloro-3-(trifluoromethyl)phenyl]-N-methyl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide</del>	
77	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
78	<del>N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[[3-(pyridin-2-ylamino)phenyl]oxy]acetamide</del>	
79	N-[2-fluoro-5-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
80	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	

Table 3		
Entry	Name	Structure
81	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea	
82	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyrimidin-5-ylphenyl)methyl]urea	
83	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
84	[3-(1H-tetrazol-1-yl)phenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
85	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	

Table 3		
Entry	Name	Structure
86	N-2-[4-chloro-3-(trifluoromethyl)phenyl]-N-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
87	2-{{4-chloro-3-(trifluoromethyl)phenyl}oxy}-N-[3-(1H-tetrazol-1-yl)phenyl]acetamide	
88	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-methyl-4-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
89	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{4-(1H-1,2,3-triazol-1-yl)phenyl}oxy}acetamide	
90	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-fluoro-4-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	

Table 3		
Entry	Name	Structure
91	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
92	<del>N-([4-chloro-3-(trifluoromethyl)phenyl]amino)carbonyl)- 3-(1H-tetrazol-1-yl)benzenesulfonamide</del>	
93	<del>N-([4-chloro-3-(trifluoromethyl)phenyl]amino)carbonyl)- N-methyl 3-(1H-tetrazol-1-yl)benzenesulfonamide</del>	
94	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2- [(4-pyridin-3-ylphenyl)oxy]acetamide	
95	2-({4-[2,4-bis(methoxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-fluoro-3-(trifluoromethyl)phenyl]acetamide	

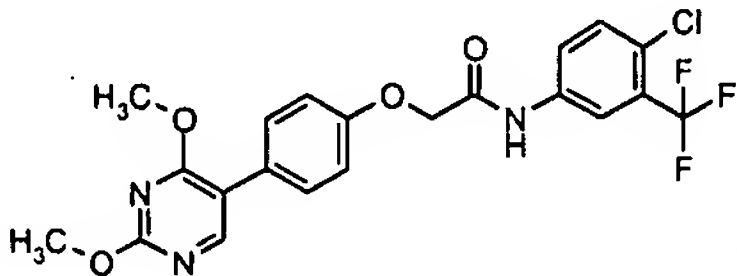
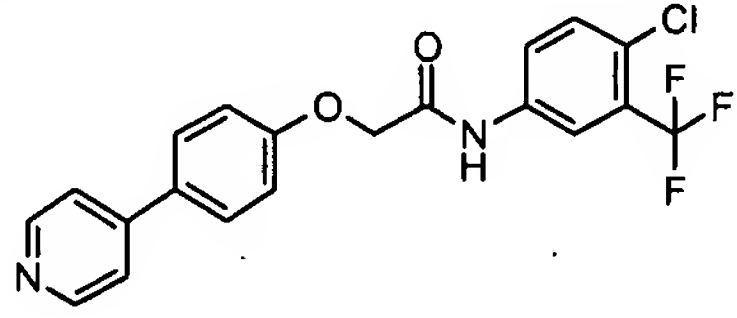
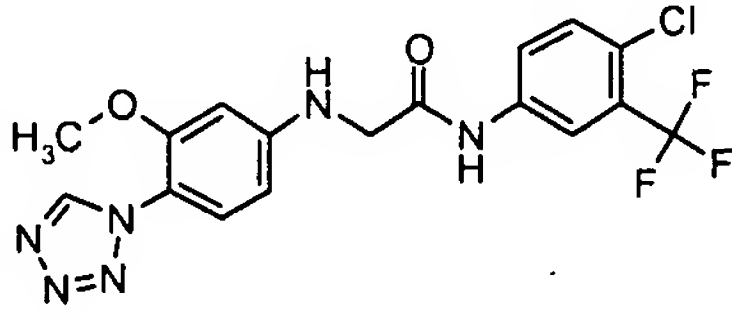
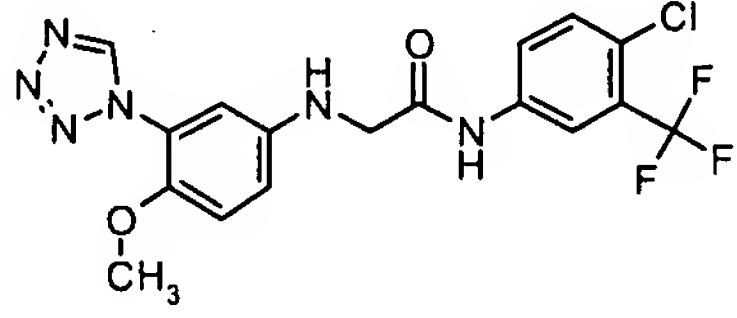
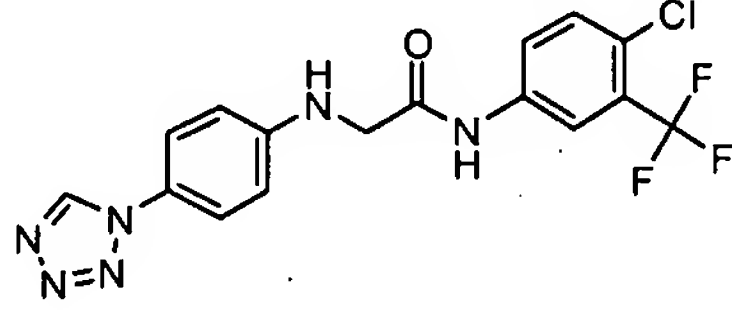
Table 3		
Entry	Name	Structure
96	2-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-chloro-3-(trifluoromethyl)phenyl]acetamide	
97	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-4-ylphenyl)oxy]acetamide	
98	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(methyloxy)-4-(1H-tetrazol-1-yl)phenyl]glycinamide	
99	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(methyloxy)-3-(1H-tetrazol-1-yl)phenyl]glycinamide	
100	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(1H-tetrazol-1-yl)phenyl]glycinamide	

Table 3		
Entry	Name	Structure
101	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(2,3,5,6-tetrafluoro-4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
102	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-tetrazol-1-yl)phenyl]methyl}urea	
103	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
104	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea	
105	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	

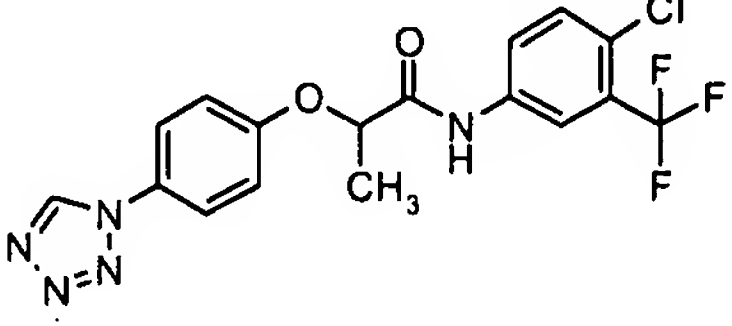
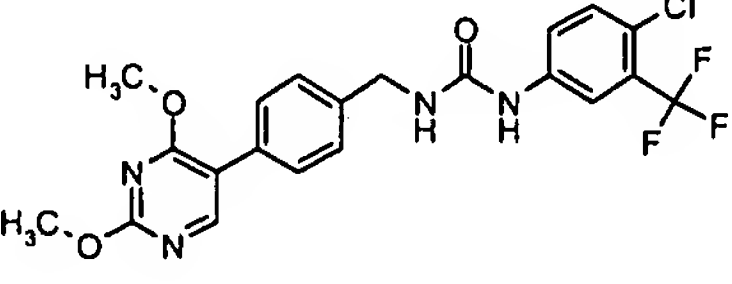
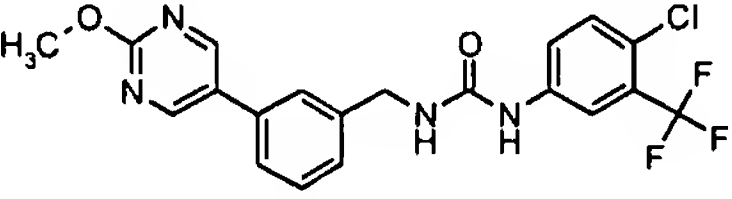
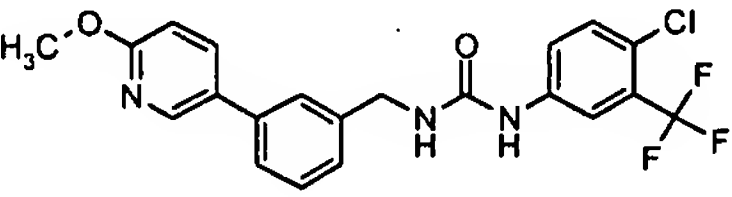
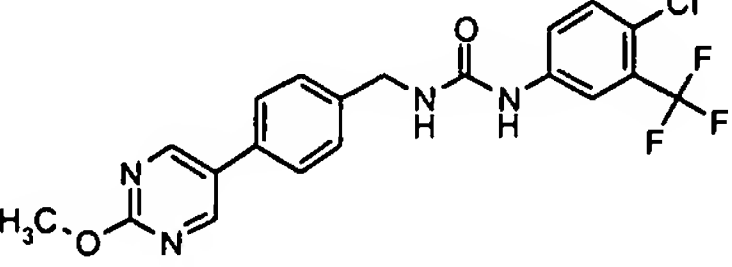
Table 3		
Entry	Name	Structure
106	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[4-(1H-tetrazol-1- yl)phenyl]oxy} propanamide	
107	N-({4-[2,4-bis(methyloxy)pyrimidin-5- yl]phenyl} methyl)-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea	
108	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({3-[2-(methyloxy)pyrimidin-5- yl]phenyl} methyl)urea	
109	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({3-[6-(methyloxy)pyridin-3- yl]phenyl} methyl)urea	
110	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({4-[2-(methyloxy)pyrimidin-5- yl]phenyl} methyl)urea	

Table 3		
Entry	Name	Structure
111	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[6-(methyloxy)pyridin-3-yl]phenyl} methyl)urea	
112	<del>1,1-dimethylethyl 2-{4-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-indole-1-carboxylate</del>	
113	<del>N-({[4-chloro-3-(trifluoromethyl)phenyl]amino} carbonyl)-4-(1H-tetrazol-1-yl)benzenesulfonamide</del>	
114	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(2H-tetrazol-5-yl)phenyl]glycinamide	
115	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2,6-difluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3		
Entry	Name	Structure
116	(3-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
117	(3-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
118	(3-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
119	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[4-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
120	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-3-ylphenyl)hydrazinecarboxamide	

Table 3		
Entry	Name	Structure
121	(4-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
122	(4-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
123	(4-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
124	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-4-ylphenyl)methyl]urea	
125	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-3-ylphenyl)hydrazinecarboxamide	

Table 3		
Entry	Name	Structure
126	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
127	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(4-pyrimidin-5-ylphenyl)methyl]urea	
128	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
129	(4-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	
130	(4-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	

Table 3		
Entry	Name	Structure
131	1-(4-pyridin-3-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
132	1-(4-pyrimidin-5-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
133	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea	
134	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea	
135	(3-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	

Table 3		
Entry	Name	Structure
136	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
137	<del>N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide</del>	
138	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
139	<del>N-[[3-(6-aminopyridin-3-yl)phenyl]methyl]-N' [4-chloro-3-(trifluoromethyl)phenyl]urea</del>	
140	<del>N-[[4-(6-aminopyridin-3-yl)phenyl]methyl]-N' [4-chloro-3-(trifluoromethyl)phenyl]urea</del>	
141	<del>N-[[3-(2-aminopyrimidin-5-yl)phenyl]methyl]-N' [4-chloro-3-(trifluoromethyl)phenyl]urea</del>	

Table 3		
Entry	Name	Structure
142	<del>N-[[4-(2-aminopyrimidin-5-yl)phenyl]methyl]-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea</del>	
143	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyridin-3-ylphenyl)ethyl]urea	
144	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyrimidin-5-ylphenyl)ethyl]urea	
145	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[4-(1H-indol-2-yl)phenyl]oxy]acetamide	
146	<del>N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(isoquinolin-7-yloxy)acetamide</del>	

Table 3		
Entry	Name	Structure
147	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-4-ylphenyl)hydrazinecarboxamide	
148	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-4-ylphenyl)hydrazinecarboxamide	
149	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-4-ylphenyl)methyl]urea	
150	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-quinoxalin-6-ylphenyl)methyl]urea	
151	<del>methyl 3-amino-6-(3-[[[4-chloro-3-(trifluoromethyl)phenyl]amino]carbonyl]amino]methylphenyl)pyrazine-2-carboxylate</del>	

Table 3		
Entry	Name	Structure
152	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-quinoxalin-6-ylphenyl)methyl]urea	
153	<del>N-[[3-(2-amino-5-methylpyridin-3-yl)phenyl]methyl]-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea</del>	
154	<del>methyl 3-amino-6-(4-[[[4-chloro-3-(trifluoromethyl)phenyl]amino]carbonyl]amino]methyl}phenyl)pyrazine-2-carboxylate</del>	
155	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-chloro-4-(methyloxy)phenyl]carbamate	
156	N-[3-chloro-4-(methyloxy)phenyl]-N'--{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea	

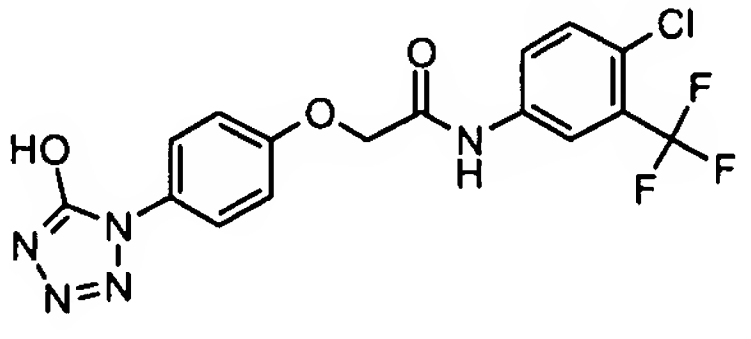
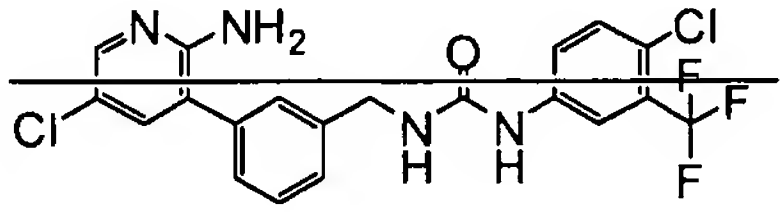
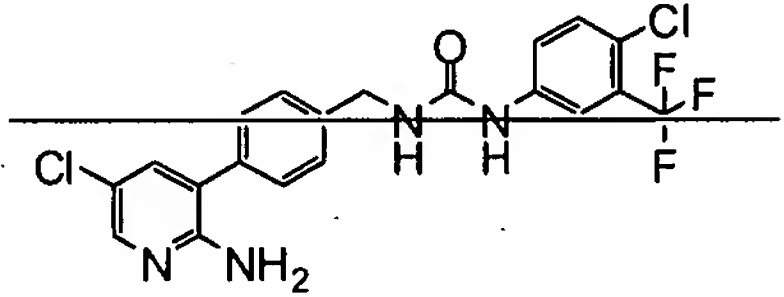
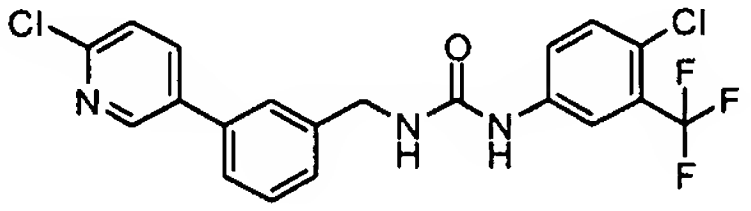
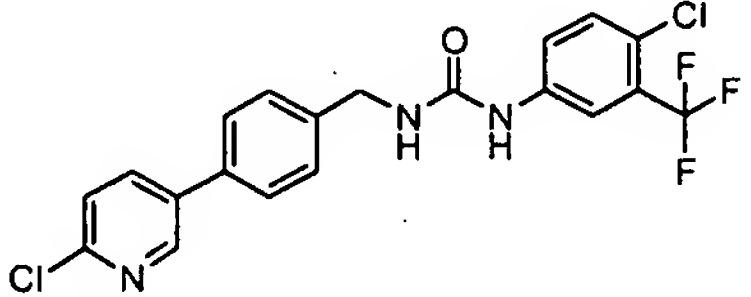
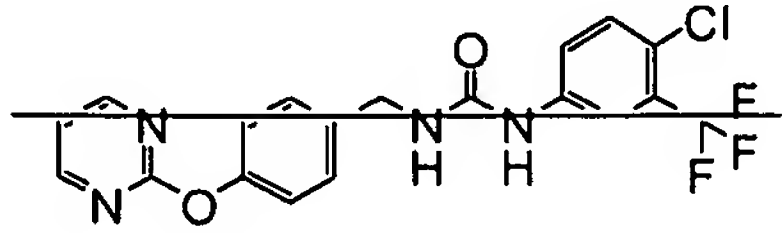
Table 3		
Entry	Name	Structure
157	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[4-(5-hydroxy-1H-tetrazol-1- yl)phenyl]oxy}acetamide	
158	<del>N-[[3-(2-amino-5-chloropyridin-3- yl)phenyl]methyl]-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea</del>	
159	<del>N-[[4-(2-amino-5-chloropyridin-3- yl)phenyl]methyl]-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea</del>	
160	N-[[3-(6-chloropyridin-3- yl)phenyl]methyl]-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea	
161	N-[[4-(6-chloropyridin-3- yl)phenyl]methyl]-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea	
162	<del>N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(pyrimidin-2- yloxy)phenyl]methyl}urea</del>	

Table 3		
Entry	Name	Structure
163	N-([4-chloro-3-(trifluoromethyl)phenyl]amino)carbonyl)-3-(1H-tetrazol-1-yl)benzamide	
164	3-amino-6-(3-([4-chloro-3-(trifluoromethyl)phenyl]amino)carbonyl)amino)methyl}phenyl)-N-[2-(dimethylamino)ethyl]pyrazine-2-carboxamide	
165	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-fluoropyridin-3-yl)phenyl]methyl}urea	
166	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-([3-[2-(methoxy)pyridin-3-yl]phenyl]methyl)urea	
167	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(6-fluoropyridin-3-yl)phenyl]methyl}urea	

Table 3		
Entry	Name	Structure
168	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({4-[2-(methoxy)pyridin-3- yl]phenyl}methyl)urea	
169	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(6-methylpyridin-3- yl)phenyl]methyl}urea	
170	<del>N-{{4-(2-amino-5-fluoropyridin-3- yl)phenyl}methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea</del>	
171	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[3-(6-methylpyridin-3- yl)phenyl]methyl}urea	
172	<del>N-{{4-(2-aminopyridin-3- yl)phenyl}methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea</del>	
173	<del>N-{{3-(2-aminopyridin-3- yl)phenyl}methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea</del>	

Table 3		
Entry	Name	Structure
174	[3-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
175	<del>[3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate</del>	
176	<del>[3-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate</del>	
177	(3-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
178	<del>N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[3-[6-(hydroxymethyl)pyridin-3-yl]phenyl]methylurea</del>	
179	N-{[3-(6-acetylpyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3		
Entry	Name	Structure
180	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[3-(6-cyanopyridin-3- yl)phenyl]methyl} urea	
181	<del>1,1 dimethylethyl (3S) 3 - ( { [ 3 amino 6 (3</del> <del>((( [ 4 chloro 3</del> <del>(trifluoromethyl)phenyl]amino} carbonyl)a</del> <del>mino]methyl}phenyl)pyrazin 2</del> <del>yl]carbonyl} amino)piperidine 1</del> <del>carboxylate</del>	
182	<del>3 amino 6 (3 - ( { ( [ 4 chloro 3</del> <del>(trifluoromethyl)phenyl]amino} carbonyl)a</del> <del>mino]methyl}phenyl) N - ( (3S) piperidin 3</del> <del>yl]pyrazine 2 carboxamide</del>	
183	<del>1,1 dimethylethyl (3S) 3 - ( { [ 3 amino 6 (4</del> <del>((( [ 4 chloro 3</del> <del>(trifluoromethyl)phenyl]amino} carbonyl)a</del> <del>mino]methyl}phenyl)pyrazin 2</del> <del>yl]carbonyl} amino)piperidine 1</del> <del>carboxylate</del>	

Table 3		
Entry	Name	Structure
184	<del>3-amino-6-(4-(((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl}phenyl)-N-[(3S)-piperidin-3-yl]pyrazine-2-carboxamide</del>	
185	[3-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
186	<del>N-[[3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl]-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea</del>	
187	<del>[6-(1H-tetrazol-1-yl)pyridin-2-yl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate</del>	
188	[3-(1H-benzimidazol-2-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3		
Entry	Name	Structure
189	<del>[3-(6-amino-2-methylpyridin-3-yl)phenyl]methyl</del> [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
190	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[5-(methylthio)pyridin-3-yl]phenyl} methyl)urea	
191	[4-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
192	<del>[4-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl</del> [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
193	<del>[4-(2-aminopyridin-3-yl)phenyl]methyl</del> [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
194	(4-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3		
Entry	Name	Structure
195	[4-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
196	[4-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
197	[3-(1H-tetrazol-1-yl)phenyl]methyl 1,3-benzothiazol-2-ylcarbamate	
198	[3-(1H-tetrazol-1-yl)phenyl]methyl (5-bromopyridin-2-yl)carbamate	
199	(3-pyridin-3-ylphenyl)methyl (3,5-dimethylphenyl)carbamate	
200	(3-pyridin-3-ylphenyl)methyl [5-chloro-2-(methoxy)phenyl]carbamate	

Table 3		
Entry	Name	Structure
201	[4-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
202	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2-(methoxy)phenyl]carbamate	
203	<del>(4-pyrimidin-5-ylphenyl)methyl (3,4-dimethylphenyl)carbamate</del>	
204	<del>(3-pyridin-3-ylphenyl)methyl (3,4-dimethylphenyl)carbamate</del>	
205	<del>1,1-dimethylethyl 3-(((3-amino-6-(3-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)pyrazin-2-yl)carbonyl)amino)piperidine-1-carboxylate</del>	

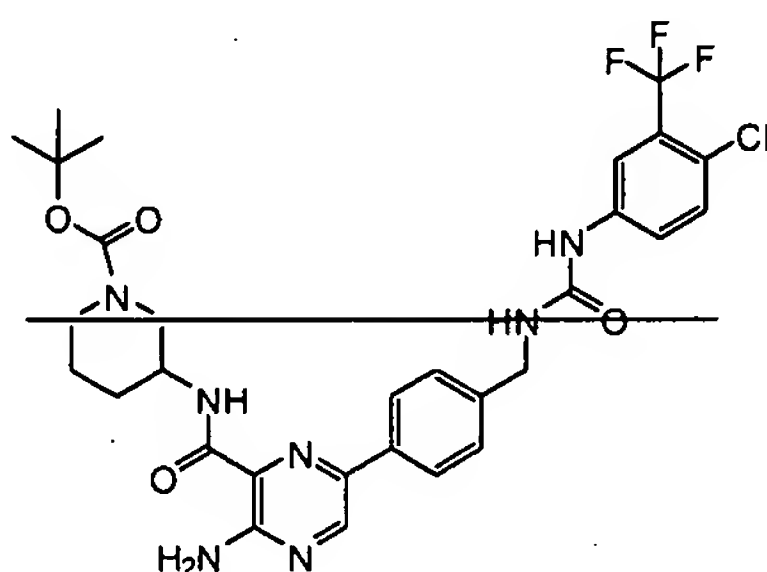
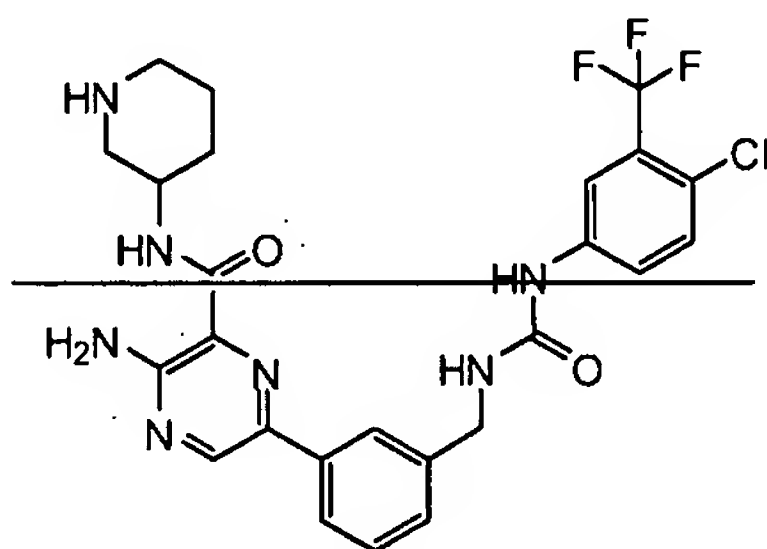
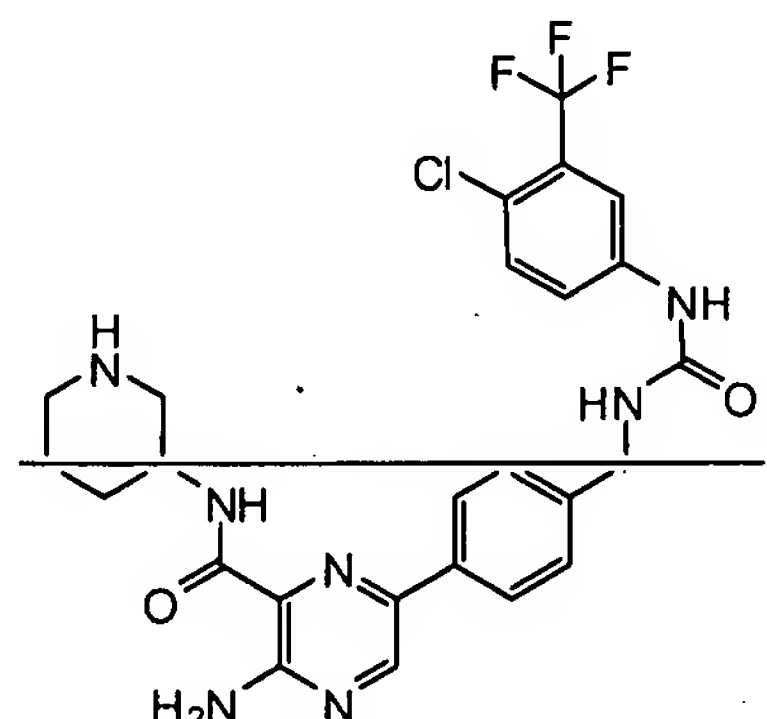
Table 3		
Entry	Name	Structure
206	<del>1,1 dimethylethyl 3 ( (3 amino 6 (4</del> <del>(((4 chloro 3</del> <del>(trifluoromethyl)phenyl)amino} carbonyl)a</del> <del>mino}methyl}phenyl)pyrazin 2</del> <del>yl}carbonyl} amino)piperidine 1</del> <del>carboxylate</del>	
207	<del>3 amino 6 (3 (((4 chloro 3</del> <del>(trifluoromethyl)phenyl)amino} carbonyl)a</del> <del>mino}methyl}phenyl) N piperidin 3</del> <del>ylpyrazine 2 carboxamide</del>	
208	<del>3 amino 6 (4 (((4 chloro 3</del> <del>(trifluoromethyl)phenyl)amino} carbonyl)a</del> <del>mino}methyl}phenyl) N piperidin 3</del> <del>ylpyrazine 2 carboxamide</del>	

Table 3		
Entry	Name	Structure
209	<del>1,1 dimethylethyl 4-((3-amino-6-(3-  ((4-chloro-3-(trifluoromethyl)phenyl)amino)-  carbonyl)amino)methyl)phenyl)pyrazin-2-  yl)carbonyl)piperazine-1-carboxylate</del>	
210	<del>1,1 dimethylethyl 4-((3-amino-6-(4-  ((4-chloro-3-(trifluoromethyl)phenyl)amino)-  carbonyl)amino)methyl)phenyl)pyrazin-2-  yl)carbonyl)piperazine-1-carboxylate</del>	
211	<del>N-((3-(5-amino-6-(piperazin-1-  yl)carbonyl)pyrazin-2-yl)phenyl)methyl)-  N'-(4-chloro-3-(trifluoromethyl)phenyl)urea</del>	

Table 3		
Entry	Name	Structure
212	<del>N-({4-[5-amino-6-(piperazin-1-ylcarbonyl)pyrazin-2-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea</del>	
213	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-pyrazol-4-yl)phenyl]methyl}urea	
214	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-pyrazol-4-yl)phenyl]methyl}urea	
215	[3-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

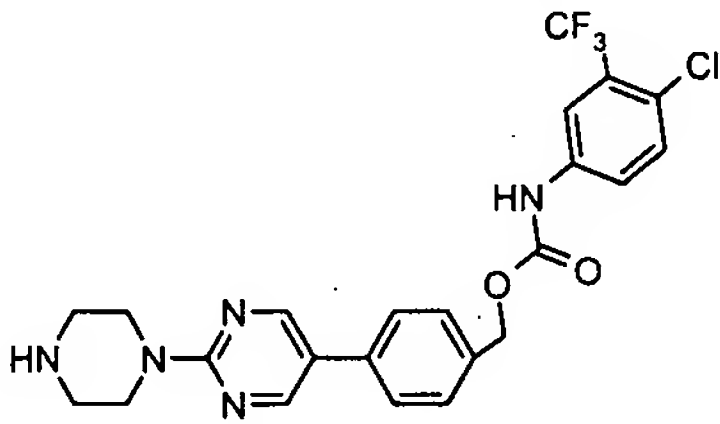
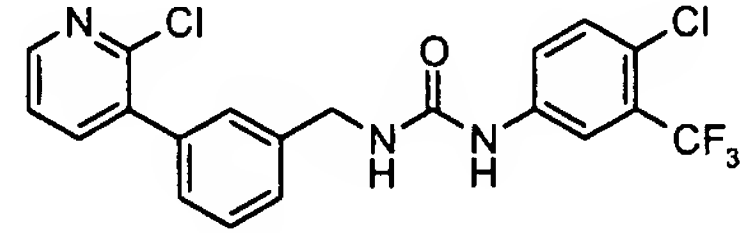
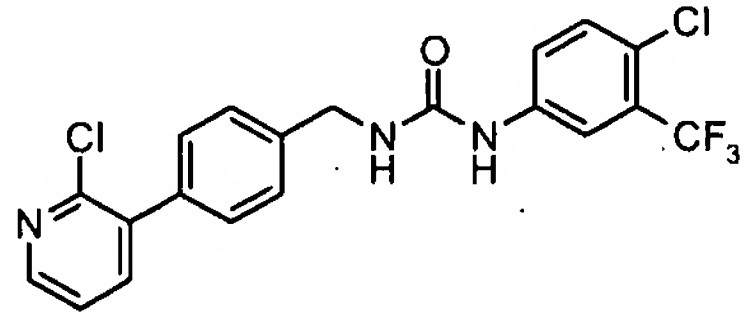
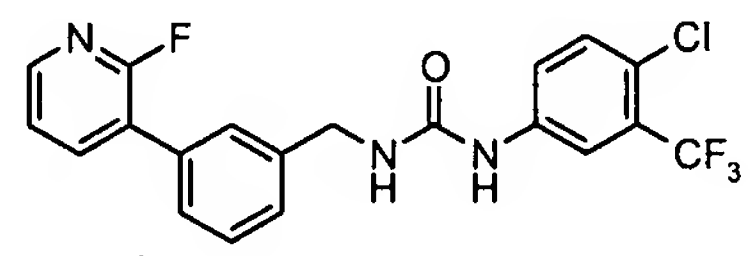
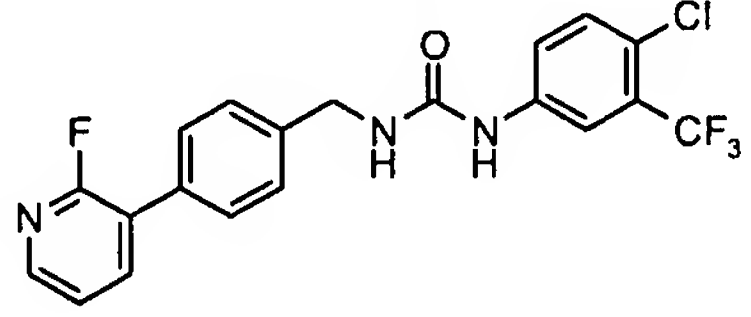
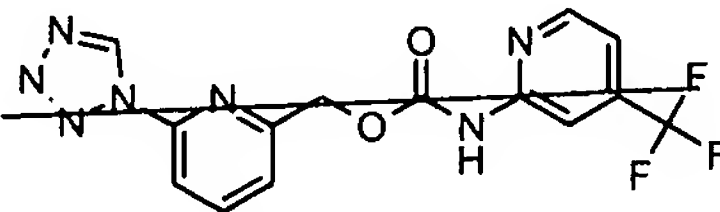
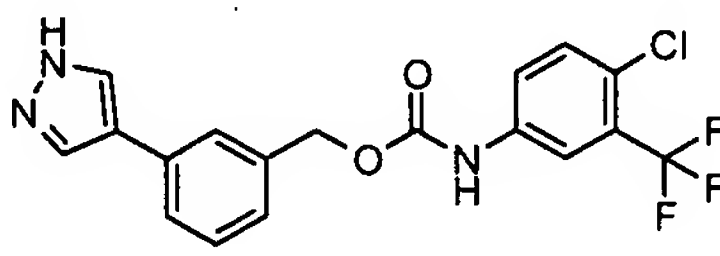
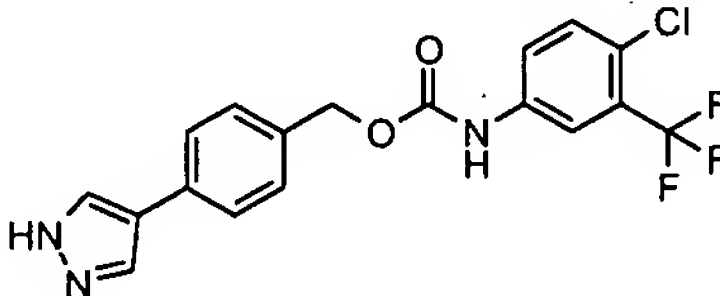
Table 3		
Entry	Name	Structure
216	[4-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
217	N- {[3-(2-chloropyridin-3-yl)phenyl]methyl} -N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
218	N- {[4-(2-chloropyridin-3-yl)phenyl]methyl} -N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
219	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[3-(2-fluoropyridin-3-yl)phenyl]methyl} urea	
220	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(2-fluoropyridin-3-yl)phenyl]methyl} urea	

Table 3		
Entry	Name	Structure
221	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-(trifluoromethyl)phenyl]carbamate	
222	[3-(1H-tetrazol-1-yl)phenyl]methyl [6-(trifluoromethyl)pyridin-2-yl]carbamate	
223	[3-(1H-tetrazol-1-yl)phenyl]methyl [4-(trifluoromethyl)pyridin-2-yl]carbamate	
224	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[5-(methylthio)pyridin-2-yl]phenyl} methyl)urea	
225	[3-(2,6-dimethylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3		
Entry	Name	Structure
226	{3-[5-(methoxy)pyridin-3-yl]phenyl}methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
227	<del>2,3'-bipyridin-6-ylmethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate</del>	
228	<del>(6-pyrimidin-5-ylpyridin-2-yl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate</del>	
229	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-isoquinolin-4-ylphenyl)methyl]urea	
230	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-isoquinolin-4-ylphenyl)methyl]urea	

Table 3		
Entry	Name	Structure
231	[6-(1H-tetrazol-1-yl)pyridin-2-yl]methyl [4-(trifluoromethyl)pyridin-2-yl]carbamate	
232	[3-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
233	[4-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

28. (previously presented) A pharmaceutical composition comprising the compound according to claim 1 and a pharmaceutically acceptable carrier.
29. (cancelled)
30. (withdrawn from consideration, currently amended) A method for modulating the *in-vivo* activity of a kinase, the method comprising administering to a subject an effective amount of the compound according to claim 1. ~~claim 1 or a compound selected from N-naphthalen-1-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-(phenyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(3,4-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-~~

~~acetamide, N-(2,3-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,4-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,5-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(3,5-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,6-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-ethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,6-diethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2-(ethyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[3-(ethyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2,4-bis(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-(dimethylamino)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,3-dichlorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-chloro-3-methylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-bromophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2-fluorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-fluorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-N-[2-(trifluoromethyl)phenyl]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-N-[3-(trifluoromethyl)phenyl]-acetamide, methyl-4-[[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetyl]amino]benzoate, ethyl-4-[[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetyl]amino]benzoate, 3-[[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetyl]amino]benzoic acid, N-[3-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy]-acetamide, N-(4-chlorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-aminophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, and N-(4-acetylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide.~~

31. (withdrawn from consideration) The method according to claim 30, wherein the kinase is c-Kit.

32. (withdrawn from consideration) The method according to claim 31, wherein modulating the *in vivo* activity of c-Kit comprises inhibition of c-Kit.

33. (withdrawn from consideration, currently amended) A method of treating rheumatoid arthritis, graft-host diseases, multiple sclerosis, psoriasis, arteriosclerosis, myocardioinfarction, ischemia, stroke, restenosis; interbowel diseases, osteoarthritis, macular degeneration, or diabetic retinopathy, ~~diseases or disorders associated with uncontrolled, abnormal, and/or unwanted cellular activities~~, the method comprising administering, to a mammal in need thereof, a therapeutically effective amount of the compound or the pharmaceutical composition as described in claim 1 ~~or a compound, or a pharmaceutical composition comprising said compound, selected from N-naphthalen-1-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-(phenyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(3,4-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,3-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,4-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,5-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(3,5-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,6-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-N-(2,4,6-trimethylphenyl)-acetamide, N-(2-ethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-ethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,6-diethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2-(ethyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[3-(ethyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2,4-bis(methyl-oxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-(dimethylamino)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,3-dichlorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-chloro-3-methylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-bromophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2-fluorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-fluorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-N-[2-(trifluoro-methyl)phenyl]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-N-[3-(trifluoromethyl)phenyl]-acetamide, methyl-4-[[3-(1H-tetrazol-1-~~

~~yl)phenyl]oxy}acetyl)amino] benzoate, ethyl 4 [([3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, 3-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(methyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-chlorophenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-aminophenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, and N-(4-acetylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide.~~

34. (withdrawn from consideration, currently amended) A method of screening for modulators of c-Kit, the method comprising combining the compound according to claim 1 ~~or a compound selected from~~ N-naphthalen-1-yl-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(phenoxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,4-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,4-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,5-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-([3-(1H-tetrazol-1-yl)phenyl]oxy}N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-ethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-diethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(ethyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2,4-bis(methyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(dimethylamino)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-chloro-3-methylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-fluorophenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-([3-(1H-

~~tetrazol-1-yl)phenyl]oxy} N-[2-(trifluoromethyl)phenyl]—acetamide, —2-[[3-(1H-tetrazol-1-yl)phenyl]oxy} N-[3-(trifluoromethyl)phenyl]—acetamide, —methyl-4-[[([3-(1H-tetrazol-1-yl)phenyl]oxy)acetyl]amino]—benzoate, —ethyl-4-[[([3-(1H-tetrazol-1-yl)phenyl]oxy)acetyl]amino]—benzoate, —3-[[([3-(1H-tetrazol-1-yl)phenyl]oxy)acetyl]amino]benzoic acid, N-[3-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, N-[4-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-[2-chloro-5-(trifluoromethyl)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-(4-chlorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, N-(4-aminophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —and—N-(4-acetylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, and at least one candidate agent and determining the effect of the candidate agent on c-Kit activity.~~

35. (withdrawn from consideration, currently amended) A method of inhibiting proliferative activity in a cell, the method comprising administering an effective amount of a composition comprising the compound according to claim 1 ~~or a compound selected from N-naphthalen-1-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, N-[4-(phenyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-(3,4-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-(2,3-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-(2,4-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-(2,5-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-(3,5-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-(2,6-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —2-[[3-(1H-tetrazol-1-yl)phenyl]—oxy} N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, N-(4-ethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, N-(2,6-diethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-[2-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-[2-(ethyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-[3-(ethyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, N-[2,4-bis(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, N-[4-(dimethylamino)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]—acetamide, —N-(2,3-~~

~~dichlorophenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-chloro-3-methylphenyl)-~~  
~~2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-bromophenyl) 2-[[3-(1H-tetrazol-1-~~  
~~yl)phenyl]oxy]-acetamide, N-(2-fluorophenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide,~~  
~~N-(4-fluorophenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, 2-[[3-(1H-tetrazol-1-~~  
~~yl)phenyl]oxy]-N-[2-(trifluoromethyl)phenyl]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-~~  
~~N-[3-(trifluoromethyl)phenyl]-acetamide, methyl 4-[[3-(1H-tetrazol-1-~~  
~~yl)phenyl]oxy]-acetyl)amino]-benzoate, ethyl 4-[[3-(1H-tetrazol-1-~~  
~~yl)phenyl]oxy]-acetyl)amino]-benzoate, 3-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetyl)amino]-~~  
~~benzoic acid, N-[3-(methoxy)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-~~  
~~(methoxy)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2-chloro-5-~~  
~~(trifluoromethyl)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-chloro-3-~~  
~~(trifluoromethyl)phenyl] 2-[[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy]-acetamide, N-(4-~~  
~~chlorophenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-aminophenyl) 2-[[3-(1H-~~  
~~tetrazol-1-yl)phenyl]oxy]-acetamide, and N-(4-acetylphenyl) 2-[[3-(1H-tetrazol-1-~~  
~~yl)phenyl]oxy]-acetamide, to a cell or a plurality of cells.~~